5 moloo74521545 mer Cethy Find attached suggested, removary revisions to broot kisk Assessment. Our approach is to provide general introduction and out the for The RA. However any calculation, conclusions and eleommendations have been deleted. I donot want to present any items at this time That many be evitical to the RA and give the appearance that They can be regotiated. Your Re-Droft should be stamped DRAFT and provided under cover letter to me, by secenter 17, if at all possible. The cover letter will introduce your prop. of a preliminary draft RA It should also recognize that the RA cannot be completed until our AI report is prepared which presents all of the into to be considered in the RA. Any questions call. Mark Bogina or me 913 551 - 7528 Thanks Claim Cuts

SYNTEX AGRIBUSINESS, INC.

DRAFT

GROUND WATER RISK ASSESSMENT

ACKGROUND be completed subsequent to preparation of the Report

(population approx. 500) in southwestern Missouri. The site is approximately 30 miles southwest of Springfield Missouri.

The Syntex facility in Verona first became the focus of environmental concerns about 1960. At that time the facility was owned and operated by Hoffman-Taff, Inc. From approximately that time until 1969, the facility produced 2,4,5-trichlorophenoxy acetic acid herbicide (2,4,5-T). In 1969, Syntex Agribusiness acquired the facility and manufactured hexachlorophene from 1970 to 1971. Both 2,4,5-T and hexachlorophene manufacture resulted in the by-product formation of dioxins, in particular, high concentrations of 2,3,7,8-tetrachlorodibenzo-p-dioxin. Concerns over dioxin contamination triggered a series of investigations at the Verona facility. Remediation of dioxin contaminated soils has been addressed in the Record of Decision signed May 5, 1988.

The Agency for Toxic Substances and Disease Registry (ATSDR) performed several toxicology consultations at the request of the U.S. EPA. These consultations (ATSDR 1989 as ammended 1990) specifically addressed the public health implications of consuming dioxin contaminated fish from the Spring River. Groundwater consumption was not addressed in these consultations.

Groundwater contamination, is being treated as a separate operable unit for the purposes of remediation. The risk assessment deals exclusively with chemicals of concern in groundwater associated with the Syntex Agribusiness, Inc. Verona facility. No dioxins or furans were detected in groundwater at the required quantitation limits.

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PHYSICAL SETTING

The major portion of the 180 acre Syntex facility is located within the 100 year flood plain of the Spring River. The facility is west of Verona and on the east bank of the Spring River. The Spring River originates south of Verona and flows northward along the western outskirts of town.

HYDROGEOLOGIC ASSESSMENT

The Syntex facility is situated in the flood plain of the Spring River Basin. The Spring River flows north in the vicinity of the facility through an alluvial valley that is underlain by limestone bedrock. The alluvial sediments consist of several feet of silty loam underlain by gravel between 10 feet and 30 feet thick. Underlying the alluvium is limestone with documented karst features. Flooding in the valley is common.

Dioxins have not been reported in the ground water in the vicinity of the Syntex facility. Aquifers that might be impacted by previous site activities include the upper alluvial aquifer and the lower limestone aquifer. Potential receptors for ground water contamination are identified in the event that contaminants from the Syntex facility have entered the ground water system.

Suspected pathways for contaminant migration in ground water to potential receptors include infiltration of contaminants from the facility downward to the ground water table during precipitation events and then transportation hydraulically downgradient in either of the aquifer systems. Alternatively, recurring flooding of the Spring River may have led to the sweeping away of contaminated sediments at the facility and redeposition of these sediments down-valley. Infiltration from precipitation could potentially transport contaminants downward to the ground water table.

In the event that contaminants from the site have entered ground water beneath the Syntex facility, potential receptors would include private well owners hydraulically downgradient of the facility or well users in which pumping of their wells could modify the local flow direction of ground water and

potentially draw in contaminants. Wells subject to potential contamination would be screened either in the valley alluvium or limestone bedrock underlying the alluvium and adjacent to the valley.

> A record search of well logs obtained from the Missouri Department of Natural Resources. Division of Geology and Land Survey (DGLS) indicate no well logs on-file for private wells within two miles of the facility that are set within alluvial sediments or limestone bedrock of the flood plain. available for private wells within two miles of the Syntex facility but most of these wells are either well outside the flood plain beyond the influence of potential contamination or hydraulically upgradient (Figure 1). Two wells are located outside of the flood plain, one approximately 1,000 feet northwest of the trench area, and the second approximately 2.500 feet north-northeast of These well locations may be indicative of potential the trench area. receptors.

> According to the U.S.G.S. 7.5 Ninute map of the Verona Quadrangle, at least nineteen houses are located down-valley of the facility and within or adjacent to the flood plain. Well logs are not on-file with the DGLS for these houses, but this does not exclude the possibility of wells being present. houses have private wells, they may be identified as potential receptors.

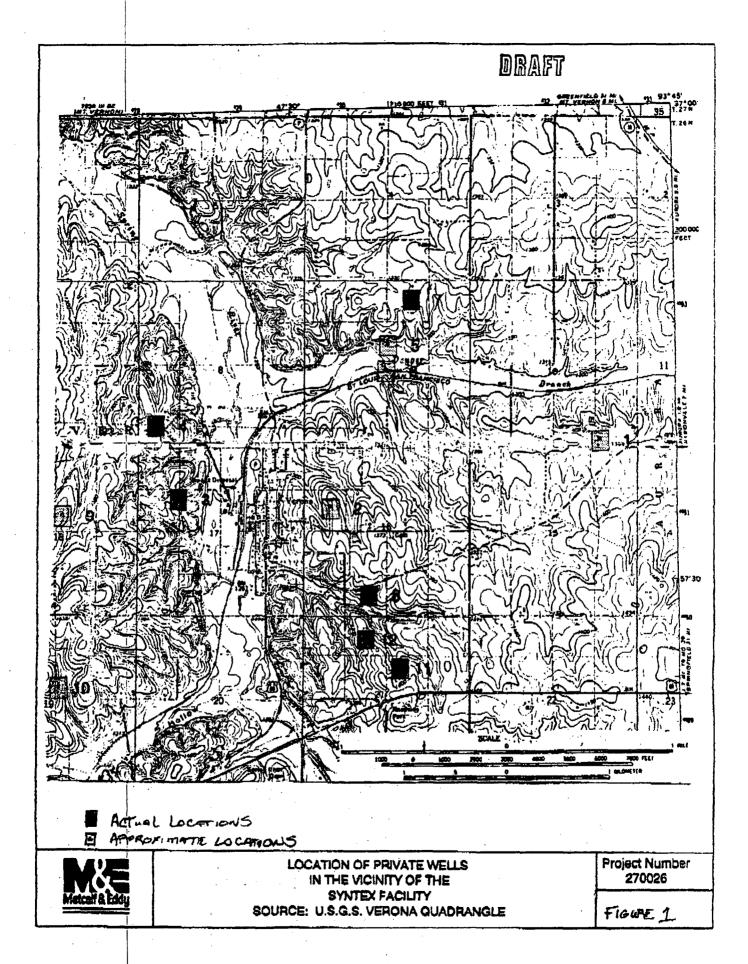
The Syntex Agribusiness acility in Vivona is considered an industrial POTENTIALLY EXPOSED POPULATIONS 5.76. Deed restrictions are in place at The site maintaining this status impersituity.

Future potentially exposed populations include existing or future downgradient residential wells and the development of future residences on the Syntex property which would include placement of onsite wells. The focus of this risk assessment is on any future residential devlopment, of the site

IDENTIFICATION OF EXPOSURE PATHWAYS

However fature development on The Syntex property is highly

Domestic use of unremediated groundwater is the only exposure pathway which will be considered in this risk assessment.



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summary of analytic results for quarterly ground water sampling performed during 1989 and 1990. Although other ground water sampling was performed in 1987, these data were not examined in the present risk assessment. The contaminants of concern presented in table 1 were selected for further study based on a validation of the ground water sampling results, comparison of sampling results to background concentrations, and whether the compounds of interest are considered to be essential nutrients or toxic. Additional

information can be considered subsequent to compilation and discussion of all available data in the RI Report.

EXPOSURE PATHWAYS AND EXPOSURE ROUTES

Exposure pathways describe the movement of chemicals from sources to locations (exposure points) where groups of people (receptors) may come in contact with the chemicals. This movement involves release of chemicals from the source to an intermediate environmental transport medium (groundwater in this case) and the receptor point. Exposure routes describe the modes of contact with and intake of contaminated media and chemicals at exposure points.

For this risk assessment, unremediated groundwater is the exposure pathway under consideration. Three potential routes of exposure will be evaluated on a chemical specific basis;

- * ingestion of chemicals in groundwater
- * dermal absorption while showering
- * inhalation of vapor phase volatiles while showering

* Irrigation

QUANTIFICATION OF EXPOSURE

Estimates of exposure levels for each contaminant are required for quantitative risk characterization. The basic equation to calculate human intake of an environmental contaminant is:

TENTATIVE

potential

TABLE 1 SUMMARY OF GROUND WATER RESULTS FOR CONSTITUENTS OF CONCERN AT THE SYNTEX FACILITY (Mo/L) *

<u> </u>		verona			
COMPOUND	MINIMUM	MUMIXAM	MEAN	STANDARD DEVIATION	UPPER LIMIT 95% C.I.
Ba	0.047	10.5	1.6822	2.2068	6.0958
Be	0.0029	0.06	0.0115	0.0149	0.0413
Cd	0.0024	0.061	0.0133	0.0196	0.0524
Cr	0.005	30.2	1.4682	5.7819	13.0356
Pb	0.001	230	7.9943	41.9554	91.9050
NI	0.04	1.9	0.1053	0.3624	0.8302
Zn	0.055	34	1.9785	7.1199	16.2182
Benzene	0.001	0.25	0.0256	0.0617	0.1491
Trichicroethene	0.005	0.25	0.0260	0.0616	0.1492
Toluene	0.005	0.25	0.0310	0.0863	0.1638
Chlorobenzene	0.002	0.25	0.0415	0.0670	0.1757
Ethyl Benzene	0.005	0.25	0.0272	0.0614	0.1500
Xylenes, total	0.001	0.25	0.0361	0.0658	0.1676
1,4-Dichierobenzene	0.01	0.11	0.0228	0.0293	0.0814
Methoxychlor	0.0005	0.0031	0.0013	0.0009	0.0031

In the final list of constituents of concern will be compited and considered subsequent to completion of the PI Report.

DI = C x HIF

where:

- DI = Daily intake (mg of chemical per kg of body weight per day)
- C = Concentration of the chemical
- HIF Human intake factor (units of medium per kg of body weight per day)

Intake variables have been selected so that the combination of all intake variables results in an estimate of a reasonable maximum exposure (RME) for the groundwater pathway. The RME is defined by US EPA (1989a) as the maximum exposure that is reasonably expected to occur at or near a given site.

METHODS FOR ESTIMATING INTAKE LEVELS

.The site

and downgrad. of the property now

It is assumed that there could be future residences on the property now occupied by the Syntex Agribusiness facility. Contaminants in soil have already been addressed in the Record of Decision For Final Management of Dioxin Contaminated Soil and Equipment at Syntex Agribusiness, Inc., Verona Missouri, April 29, 1988. This risk assessment will address potential future exposures to groundwater contaminants at the Syntex Agribusiness site. In this The exposure scenario, it is assumed that persons will reside on the site and use unremediated groundwater for all domestic purposes including drinking and cooking.

Three specific routes of exposure to contaminants in ground water are addressed: ingestion, dermal absorption while showering, and inhalation of volatile components while showering. By convention, daily intakes of the contaminants of concern will be estimated for an average 70 kg adult and for a 16 kg child (1 to 6 years old). It is assumed that the adult will reside on the Syntex site for a maximum of 30 years (national 90th percentile time at one residence as estimated by U.S. EPA). During this 30 years, it is assumed that exposure to contaminants in groundwater occurs 365 days per year. Children are assumed to be exposed for five years. Subchronic intake for

same

adults and children will be based on five years of exposure. Chronic intakes will be calculated only for adults based on 30 years of residence on the site. Carcinogenic effects will be estimated for 30 year exposure duration for adults and 5 years for children, both averaged over a 70 year lifetime.

INGESTION OF GROUNDWATER

Daily intake of contaminants of concern through ingestion of groundwater will be estimated for both an average adult and child. A 70 kg adult is assumed to drink 2 liters of water or beverages made with groundwater every day, 365 days per year for either 5 years (subchronic effects) or 30 years (chronic noncarcinogenic and/or carcinogenic effects). A 16 kg child is assumed to drink 1 liter of water or beverages every day for 365 days per year for 5 years. It is further assumed that the administered dose, as estimated by the daily intake, is also the absorbed dose

DERMAL ABSORPTION WHILE SHOWERING

It is assumed that any future residents on the Syntex Agribusiness site will shower every day for 365 days per year using unremediated ground water. Dermal absorption of semi-volatile or inorganic compounds may occur during showering. A 70 kg adult is assumed to have a skin surface area of 18,150 cm². A 16 kg child is assumed to have a corresponding skin surface area of 9,400 cm². Shower exposures are assumed to last a maximum of 15 minutes. Where available, chemical specific permeability constants will be used to estimate the amount of the chemical absorbed through the skin and into the peripheral circulation. Where chemical specific permeability constants are not available, standard values for soluble and insoluble compounds will be utilized.

As/described for the ingestion route of exposure, showering will be assumed to occur 365 days/year for 30 years for adults and 365 days/year for 5 years for children. Subchronic, chronic noncarcinogenic, and carcinogenic effects will be estimated as described previously.



INHALATION OF VAPOR PHASE CHEMICALS WHILE SHOWERING

Under shower conditions in a confined area, volatile organic chemicals present in groundwater would be expected to enter vapor phase. Consequently, inhalation of volatile organic compounds while showering could contribute to an individual's total exposure. The amount of each volatile compound vaporizing into the air during a 15 minute shower will be estimated based on the concentration in ground water. The enclosed volume of a 1 m x 1 m x 2.5 m shower stall is assumed to be 2.5 m³. Determination of chemical concentrations of waterborne VOCs in shower stall air were determined by using a flow rate of 15 gallops/minute (56.8 l/minute), multiplied by a shower duration of 12 minutes (U.S. EPA, 1989a), divided by shower stall dimensions of 2.5 m³. The resulting quantity of water per unit volume from which VOCs could volatilize is 273 l/m³.

While showering, the individual is assumed to inhale the estimated mean concentration of the chemical for 0.25 hours each day for 365 days per year. Respiration rate is assumed to be 0.5 m³/hour for both adults and children (U.S. EPA 1989). All other assumptions are as described previously.



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Ingestion of Groundwater

Equation (U.S.EPA, 1989a):

Intake (mg/kg day) = $\frac{CW \times IR \times EF \times ED}{BW \times AT}$

/ARIABLI	E VARIABLE DESCRIPTION	GROUNDWATER ASSUMPTIONS
CW	Chemical Concentration in Water (mg/liter)	Constituent Concentration in Groundwater
IR	Ingestion Rate (liters/day)	<pre>2 liters/day (adult) 1 liter/day (child)</pre>
EF	Exposure Frequency (days/year)	365 days/year
ED	Exposure Duration (years)	30 years for adults (chronic and carcinogenic effects)
		5 years for adults (subchronic effects)
		5 years for child (subchronic and carcinogenic effects)
BW	Body Weight (kg)	70 kg (adult) 16 kg (child)
AT	- Averaging Lime (period over which expesure is averaged adays)	25,550 days (carcinogenic effects)
		10,950 days (chronic noncarcinogenic effects)
		1,825 days (subchronic effects)

TABLE 5 -

Derma Contact with Groundwater

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Equation (U.S.EPA, 1989a):

Intake (mg/kg day) = <u>CW x SA x PC x ET x EF x ED x CF</u>
<u>BW x AT</u>

VARIABLE	VARIABLE DESCRIPTION	GROUNDWATER ASSUMPTIONS
CM	Chemical Concentration in Water (mg/liter)	Constituent Concentration in Groundwater
SA	Skin Surface Area Available for Contact (cm)	18,500 cm ² (adult) 9,400 cm ² (child)
PC	Chemical Specific Dermal Permeability Constant (cm/hr)	Chemical Specific (cm/hr)
ET	Exposure Time	0.25 hours/day
EF	Exposure Frequency (days/year)	365 days/year
CO	Exposure Duration (years)	30 years for adults (chronic and carcinogenic effects)
		<pre>5 years for adults (subchronic effects)</pre>
		5 years for child (subchronic and carcinogenic effects)
CF	Volumetric Conversion Factor for Water (1.1) ter/1000 cm)	1 liter/1000 cm ³
BW	Body Weight (kg)	70 kg (adult) 16 kg (child)
AT	Averaging Time # (periodeover which exposure is averaged ==days)	25,550 days (carcinogenic effects)
		10,950 days (chronic noncarcinogenic effects)
		1,825 days (subchronic effects)



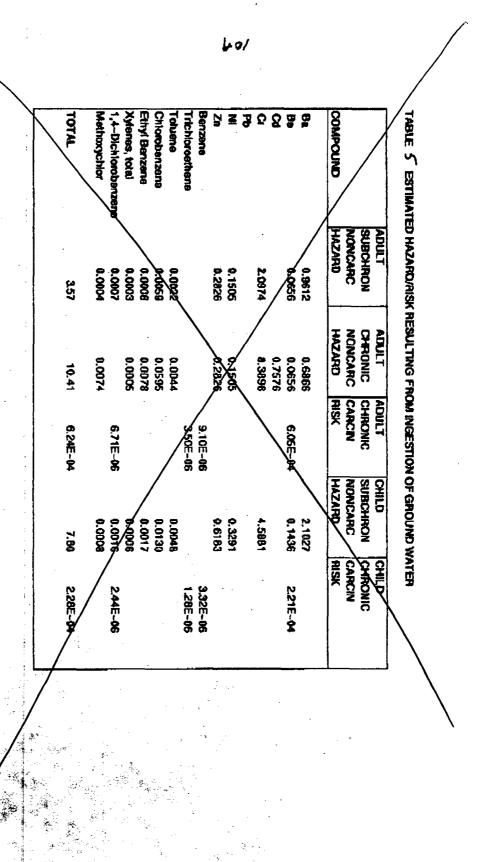
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Inhalation of Vapor Phase Chemicals While Showering with Groundwater

Equation (U.S.EPA, 1989a):

Intake (mg/kg day) = $CA \times IR \times WV \times ET \times EF \times ED \times CF$ BW x AT

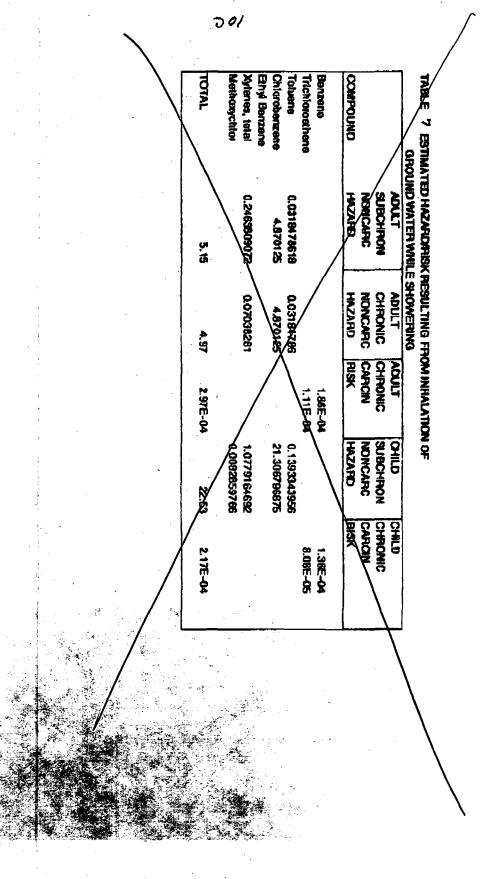
VARIABLE	VARIABLE DESCRIPTION	GROUNDWATER ASSUMPTIONS
CA	Chemical Concentration in Air During Shower (mg/m³)	Constituent Concentration in Air During Shower
WV	Water Quantity per unit volume (L/m³)	273
IR	Inhalation Rate (m³/hr)	0.6m³/hr
ET	Exposure Time (hours/day)	0.25 hours/day
EF	Exposure Frequency (days/year)	365 days/year
ED	Exposure Duration (years)	30 years for adults (chronic and carcinogenic effects)
		5 years for adults (subchronic effects)
		5 years for child (subchronic and carcinogenic effects)
BW	Body, Weight (kg)	70 kg (adult) 16 kg (child)
AT#	Averaging Time (period over which exposure is averaged - days)	25,550 days (carcinogenic effects)
		10,950 days (chronic noncarcinogenic effects)
×		1,825 days (subchronic effects)



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TABLE 6	ESTINATED HAZARD/RISK RESULTING FROM DERMAL CONTACT WITH
	GROUND WATER WHILE SHOWERING

COMPOUND	ADULT SUBCHRON NONCARC HAZARD	ADULT CHRONIC NONCARC HAZARD	ADULT CHRONIC CARCIN RISK	CHILD SUBCHRON NONCARC HAZARD	CHEO CHRONIC CARCIN RISK
	2000	0.0000		0.0744	
Ba	8.0333	0.0238		0.0741	
Be	0.0023	0.0023	2.18E-06	0.0051	7.77E-08
Cd	`	0.0375			
Cr	0.0146	0.0582		0.0323	
Pb		\ /			
Ni	0.0052	0,9052		0.0116	
Zn ·	0.0033	0.0033		0.0073	
Benzene			8,34E-08		8.66E-07
Trichiproethene	*		8.105-07		3.00E-07
Toluene	0.0005	0.0010		0.0011	•
Chlorebenzene	0,9646	0.0458		0.0102	
Ethyl Benzene	0.0002	0.0022		9,0005	
Xylenas, total	0.0001	0.0001		0.0001	, ·
1,4-Dichlorobenzane	0.0002		1.55E-06	0.0004	5.75E-07
Methoxychlor	0.0001	0.0017		0.0002	
TOTAL.	0.06	0.18	2.57E-05	0.14	9.51E-06



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EXPOSURE CONCENTRATIONS

Risk assessment maintains The conservation approach, onsidering In keeping with the conservative fature of this risk assessment; exposure point concentrations in groundwater were based on the maximum detected concentration of each chemical. /Estimations of exposure concentrations assume that levels of each chemical will remain constant over the duration of the exposure period. This is at extremely conservative assumption, given that groundwater flows would be expected to cause the concentrations of most source related contaminants to decrease over time.

I may be

TOXICITY OF CHEMICALS OF POTENTIAL CONCERN

The most critical effects of the chemicals of concern are summarized in the following tables. Brief toxicological profiles can be found in Appendix B.

RISK CHARACTERIZATION

Evaluation of Noncarcinogenic Risks:

The risk of adverse noncarcinogenic effects from chemical exposure is expressed in terms of the Hazard Index (HI). The HI is the ratio of the estimated dose which a hypothetical receptor would receive to the estimated dose level believed to be without measureable adverse effects. The estimated "safe" dose level is termed the Reference Dose (RfD). The Hazard Index is calculated as follows:

HI - CDI/RfD

HI = Hazard Index for Chronic Exposure

CDI - Chronic Daily Intake

RfD - Reference Dose

If the Hazard Index is less than 1.0, it is believed that the risk of noncarcinogenic effects is very low. If the HI exceeds 1.0, there is a risk of adverse noncarcinogenic effects. However, this risk is still likely to be

TABLE S
SYNTEX AGRIBUSINESS FACILITY, VERONA, MISSOURI
TOXIC EFFECTS ASSOCIATED WITH CHEMICALS OF CONCERN

	NONCARCINOGENIC E	FFECTS	CARCINOGENIC E	FFECTS*
CHEMICAL	ORAL	INHALATION	ORAL	INHALATION
Bartum	incr. Blood	Fetotokielty	Inecretuaive	Inconclusive
	Pressure		(Group D)	(Group D)
Seryllium .	ND	None Observed	Total Tumore (Group B2)	Lung Tumore (Group B2)
Cadmium	Renal Toxicity	Cencer	NA	Respiratory Tumors (Group B2)
Chromium III	Hepatotodolty	NA	Inconclusive (Group D)	(Group D)
Chromium VI	ND	Cancer '	NA	Lung Tumore (Group A)
	* i			
Lead	Central Nervous	Central Nervous	NA	NA
	Toxicity	Toxicity	(Group 82)	(Group BZ)
Nickel	Decr. Body Weight	Cancer	NA	Respiratory Tumors (Group A)
Zinc	Anemia	NA	inconclusive (Group D)	Inconclusive (Group D)
Ethylbenzene	Hepatorodoity and Nephrotoxicity	NA .	(Group D)	inconclueive (Group D)
Toluene	Central Nervous System Toxicity	Central Nervous System Toxicity Eye and Nose	inconclusive (Group D)	Inconclusive (Group D)
Xylenes	Hyperactivity Weight Lees	Central Narrous System Toxicity;	incombinativo (Group D)	inconclusive (Group D)
Chlorobenzene	Hepatotoxicity and Nephrotoxicity	Nose and Throat Hepatotoxicity and Nephrotoxicity	Inconclusive (Group D)	inconclusive (Group D)
Trictiloroethane		NA ·	Liver Turnors (Group B2)	Lung Tumore (Group B£)
1,4-Diahlorobenzene		Hepatoexicity and Nephrotoxicity	Liver Tumors (Group B2)	NA
Methoxychlor	Fetotoxicity.	NA	Inconclusive (Group D)	Inconclusive (Group D)

^{*}U.S. EPA Weight of Evidence Classification:

11000

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TABLE Certifolen
Group A, Human Carcinogen
Group B, Probable Human Carcinogen
Group D, Nex Classifiable
Group E, Evidence of Noncarcinogenicity

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small since RfDs are developed in a highly conservative manner. The available oral Reference Doses for chemicals of concern are listed in Table? IN Appendix B Tables 9 10 and 1/ propert the hexaeds associated with the chemicals of potential concern.

Evaluation of Carcinogenic Risks:

Chemicals for which there is human and/or laboratory evidence of carcinogenicity will be assessed for potential lifetime cancer risk. The risk of cancer from exposure to a specific chemical under specified exposure circumstances is expressed as a probability value. This is termed the excess lifetime risk because it is in addition to an individual's already existant chances of developing cancer during his or her lifetime. The excess lifetime risk for a given chemical is calculated by multiplying the Chronic Daily Intake by the Cancer Potency Factor, also termed the Slope Factor as derived from laboratory data.

Cancer Risk = CDI x SF

Risk estimates are presented as excess cancer risk per unit of exposed population. This risk assessment will use the Slope Factors which have been derived by U.S. EPA. These calculations are extremely conservative and provide a 95% upper bound estimate of the risk. This means that the true risk may be lower than the calculated probability, but is not likely to be higher.

Because an individual may be exposed to many different potentially carcinogenic chemicals at the same time, the risk estimates are added together for all chemicals in a specific pathway and for all pathways. In this case, risk estimates will be added for all routes of exposure encompassed by the groundwater pathway. Tables 7, 10 months the summation of the individual carcinogenic risks for each potentially carcinogenic chemical of concern.

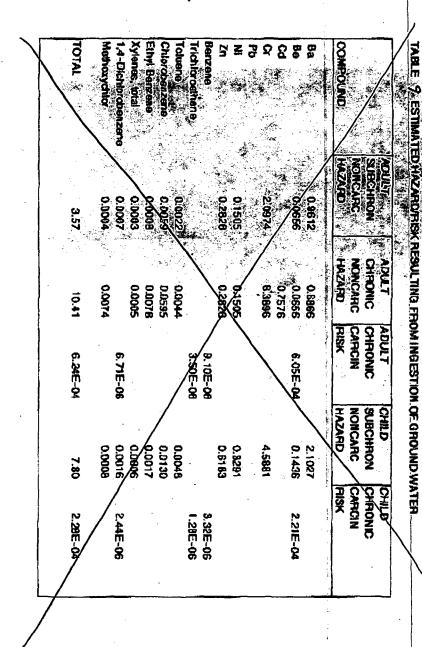
UNCERTAINTIES IN RISK ASSESSMENT

The assumptions and accompanying uncertainties associated with estimating the risks from consumption of unremediated groundwater at the Syntex Verona facility are summarized below.

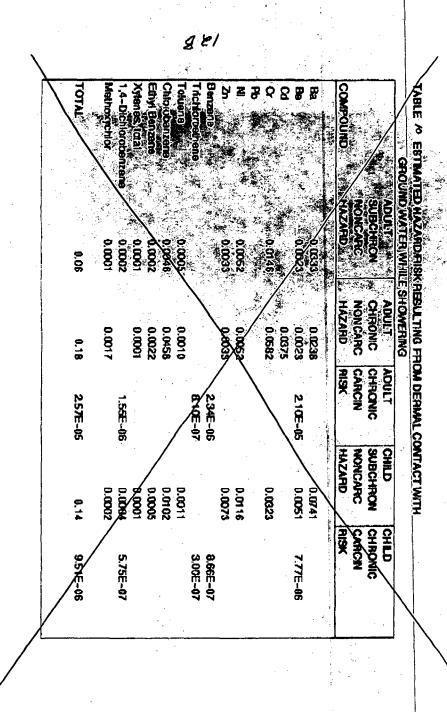
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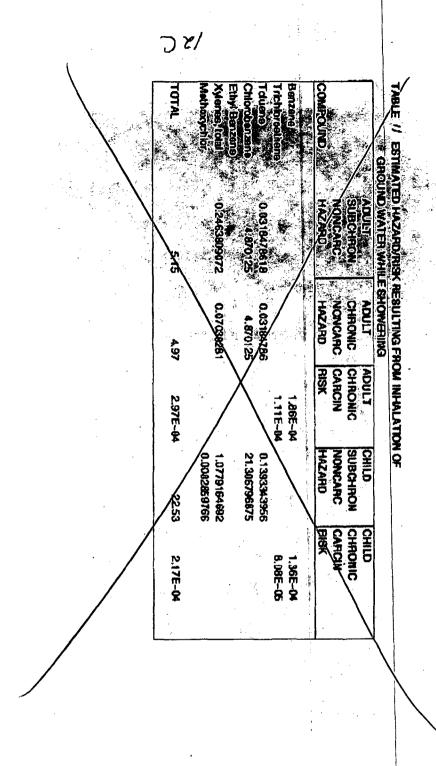
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future use questions - well not be residential

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- It is assumed that chemical concentrations will remain constant over the duration of an individual's exposure. Movement of groundwater under the Systex property is likely to produce a progressive dilution of site related contaminants, particularly after the sources have been remediated. By assuming that the measured concentrations of chemicals of concern will remain constant for the next 30 years, it is very likely that exposures will be over-estimated.
- It is conservatively assumed that the adverse effects of exposure to mixtures of chemicals will be additive (US EPA 1989a). This may or may not be true. Furthermore, humans display a diversity of responses to xenobiotic chemicals which cannot be fully incorporated into the risk assessment process.
- Extrapolation of effects observed in laboratory animals to human populations is inherently uncertain. To account for this uncertainty, a factor of 10 (or some multiple) may be used in extrapolating doses from laboratory animals to humans. In general, this is believed to increase the conservatism of the estimate
- Many of the chemicals present in groundwater associated with the Syntex Agribusiness site are frequently encountered in groundwater and in other media. Site attributable risk can only be determined to the extent that background has been adequately characterized.

There is considerable debate and uncertainty regarding the calculation of Slope Factors for carcinogenic chemicals. In order to be conservative, the U.S. EPA colculates the upper 95% confidence limit of the slope at low doses. Consequently, the actual slope factors may be lower about are very unlikely to be higher. These calculations also assume that there is no threshold for carcinogenic effects. Again, this may or may not be true for the chemicals under consideration.

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The results of any risk assessment must be interpreted in light of the inherent uncertainties. Case specific uncertainties should also be weighed during the interpretation.

CONCLUSIONS

The results of the estimation of noncancer hazards associated with the concentrations of the chemicals of concern indicate that excess hazards (greater than the value of unity accepted by U.S. EPA) may be associated with the ingestion of chromium levels and inhalation (during showering) of chlorobenzene levels found in ground water at the Syntex Pacility. Dermal contact (during showering) with the constituent concentrations in ground water does not appear to present a noncarcinogenic health hazard.

The potential cancer risk estimates associated with the concentrations of the chemicals of concern indicate that risks above the level of 1.0E-6 (level considered acceptable by U.S.EPA) may be associated with the ingestion and dermal contact (during showering) with peryl/lum levels and inhalation (during showering) of benzene and trichloroethylene levels found in ground water at the Syntex Facility.

These findings indicate that adverse human health consequences may potentially exist for receptors utilizing ground water in the vicinity of the Syntex facility.

RECOMMENDATIONS

- A detailed survey of local private wells in the vicinity of the Syntex Facility should be performed to ascertain the actual number of human receptors who may be using ground water downgradient of the site.
- The use designation of local private wells in the vicinity should be determined to demonstrate whether the ground water from the local wells is actually being utilized downstream of the site. The use designation should

TABLE 12.
SUMMARY OF HAZARDS/RISKS ASSOCIATED WITH CHEMICAL CONCENTRATIONS IN GROUND WATER

	ADULT				CHILD	
PROUND WATER	SUBCHRON NONCARC HAZARD		CHRONIC NONCARC HAZARD	CARCINOG RISK	SUBCHRONIC NONCARC HAZARD	Carcinog Risk
NGESTION	\searrow	3.57	10.41	6.24E-04	7.80	2.28E-04
SHOWERING/DERMAL CON		2.06	0.18	2.57E-05	0.14	9.51E-0
BHOWERING/INHALATION	/ !	5.75	4.97	2.97E-04	22.53	2.17E-0
TOTAL		3.78-	15.57	8.47E-04	\$0.48	4:54E-0
*****	international service recommend	4000,04004.0				

APPENDIX A

Ground Water Sampling Results and Selection of Constituents of Concern

TABLE A-1 GROUND WATER SAMPLING RESULTS FOR THE SYNTEX FACILITY (MGA.)

	MW1			MWS			MW6			MW7	
	SYN	SYN	EPA	SYN	SYN	EPA	SYN	SYN	EPA	SYN	SYN
COMPOUNDS	1/30/90	4/26/90	7/26/90	1/30/90	4/26/90	7/26/90	1/30/90	4/26/90	7/26/90	1/30/90	4/26/94
Al 1 mg .			_ 30	ľ	ļ	120		į.	500	1	1
As	0.01	0.01		0.01	0.01	0.049	0.015	0.022	0.072	0.016	0.01
Ba	10.5	0.296	1.5	5.42	2.76	3.4	3.26	2.29	6.1	1.9	1.1
89			0.005			0.014			0.05	}	}
Cd						0.0094	*		0.061		
Ca	119	48.5	59	97.1	71	75	149	114	190	121	67.
Q ₇	30.2	0.731	6.9	0.157	0.119	0.22	0.021	0.005	0.7	0.007	0.03
Co			0.1			0.092			0.3	5.55	
Q ₁		1	0.065			0.09		1	0.38	1	1
Fe	301	13.3	80	167	108	150	106	85.4	B10	49.2	35.
Pb	0.322	0.006	0.029	0.173	0.072	230	0.037	0.003	0.73	0.013	0.01
Mg	15.6	2.41	5.3	15.6	11.2	17	11.3	6.23	35	11.1	6.
Mn	80.8	2.41	13	28.9	11.6	13	125	8.78	81	5.87	4.1
N			1.9			0.23			0.86		
K		1	5.4	}	ł	20	}		53		1
Se	0.005	0.005	0.05	0.005	0.005	0.05	0.005	0.005		0.005	0.00
Ag	1		0.036		}	0.028			0.082		
Na	7.7	6.75	8.6	42.5	43.8	0.05	117	104	87	60.2	40.
TI			0.1						1	1	1
v	i		0.13		Ì	0.23	1		0.77		1
Zn			0.25			0.53		1	2.7		
Methylene Chloride	0.005	0.001	0.005	0.005	0.04	0.005	0.005	0.024	0.005	0.005	0.00
Benzene	0.005	0.005	0.005	0.005	0.017	0.005	0.001	0.005	0.005	0.005	0.00
Trichioroethene	0.005	0.005	0.005	0.005	0.017	0.005	0.005	0,005	0.005	0.013	0.00
Taluene	0.005	0.005	0.005	0.005	0.017	9.005	0.168	0.005	0.005	0.006	0.00
Chlorobenzene	0.005	0.005	0.005	0.064	0.017	0.029	0.169	0.016	0.005	0.02	0.01
Elhyl Benzene	0.005	0.005	0.005	0.405	0.017	0.005	0.029	0.018	0.015	0.005	0.00
Acetene	0.059	0.224	2	0.013	0.836	0.59	0.124	0.151	0.63	0.051	0.20
Carbon Disulfide	0.005	0.005	0.005	0.005	0.017	0.005	0.005	0.005	0.006	0.006	0.00
Xylenes, total	0.005	0.005	0.12	0.005	0.017	0.009	0.142	0.026	0.025	0.005	0.00
1,4-Dichlorobenzene			0.01			0.016		*	0.11		
1,2,4-Trictflorobenzane			0.01			0.01	1		0.011		
Bis(2-Ethylhekyl)Phthalate		1	0.01)	İ	0.01	[1	0.01]
Methoxychior				j							

PAREC

	MW15A		WW158					MM16			
90.	╝		Ad3	SYN	BPA	NAS	BPA	╛		€ PA	NAS
COMPOUNDS	1/31/90	4/26/90	1/30/90	1/31/90	2101/90	4/26/90	7/25/90	1/31/90	1/31/90	2/01/90	4/26/90
≥							0.2	£			
26	0.01	0.01	•	0.017		0.011		<u>9</u> .01	<u>8</u>		0.01
	0.817	2,33	0.047	1.B94		22	=	0.12	0.327		0.198
Be							0.005	0.005			
8							0.006	0.0024			
Ca	67.8	85.9	110	132		116	74	9	171		89.5
Q	0.023	0.182		0.013		200	0.Q	0.02	0.034		6,00,0
8				•			0. 8 6	0.0091			
S			-				0.025	0.025			
Fe	18.9	146		56,1		80.9	双	13	21.7		7.49
P	0.014	0.128		0.022		0,05	0.003	0.015	0.018		0.005
PAG	5.73	11.6	8.8	=======================================		10.3	6.5	19	18.5		17.4
Ma	12.9	23.7	0.22	<u>5</u> 5		77.1	4.5	0.7	287		11.3
2	-				-		0,04	0.04			
) X							સ્ટ	<u>.</u>			
, a	0.000	0.000		0,000		CUU.	0.00	0.000	0.000		0.000
8			1				9.9	0.01			<u> </u>
Za	39.3	27.8	7.4	159		95.6	16	9.6	32.7		48.6
3								0.01			
<							0.08	0.028			
27			0.088				¥	0.2 1			
Methylone Chlaride	0.005	0.28		0.01	0.25	0.667	0.005		0.125	0.005	0.109
Benzene	0.005	0.005		0.01	0.25	0.25	0.005		0.125	0.005	0.034
Trichlaroethene	0.005	0.005		0.01	20,23	0.25	0.005		0.125	0.005	0.034
Toluene	0.005	0.005		0.016	0.25	0,25	0.005		0,125	0.005	0.034
Chlorabenzene	0.005	0.005		0.11	0.25	0.25	0.005		0.125	0.005	0.034
Ethyl Benzene	0.005	0.005		0.008	0.25	0.25	0.005		0.125	0.005	0.034
Acetone	0.114	3.93		0,4	9.6	11.79	0.098		4.4	0.19	1.54
Carbon Disultide	0.005	0.005		0.01	0.25	0.25	0.005		0.125	0.005	0.034
Xylanes, total	0.005	0.005		0.032	0.25	0.25	0.021		0.125	0.006	9.034
1,4-Dichbrobenzene					0.01		0.046			0.01	
1,2,4-Trichtorobenzene					0.01		0.01			0.01	
Bis(2-Ethylhexyl)Phthalate					0.015		0. 2.			0.003	
Methoxychiar					0.0015					0.00087	

TAARO

	MWO		L F. WILL		AND THE			
	MAS	NAS	EPA	EPA	EPA	KYS	EPA	MAS
COMPOUNDS	ဗ	₩25/90	1/31/90	2/01/90	1/30/90	1/31/90	98/K	
>			8.7		N3 N3			
\$	0.09	0.019	0.01		0.014	0.01		0.01
	0.705	1.36	0.18			0.368		0.05
130			0.0029	•	,			
8			0.005	•				
Ca	74.9	99.2	130		- - - - - - - -	231		102
Cr	0.041	0.06	0.031			0.032		0.006
Co			0.024		0.012			
C.			0.043					
Fe	50.7	113	. 22		13	16.6		0.236
3	0.009	0.057	0.065		0.001	0.029		0.003
Ng	5.55	10.4	*		8.4	121	-	8.72
75	5.54	B.53	1.8		11	3.41		0.204
2			0.069		0.046			
*			7.3	•	7.3			
Sa	0.005	0.005	0.005			0.005		0.005
20			0.01					
Np	79.2	85.5	6.7		110	14.		5.97
1			0.01					
<			0.032					
Zn			0.34		0.085			
Methylese Chloride	9000	0.003		0.005	0.038	0.005	0.005	
Benzene	9000	0.005		0.006	0.005	0.005	0.005	
Trichloroethene	200.0	0.005		0.005	0,005	0.005	0.005	
Toluene	9000	0.005		0.005	0.005	0.005	0.005	
Chlorobenzene	0.049	0.087		0.005	0.002	0.002	0.006	
Ethyl Benzens	0.005	0.005		0.005	0.005	0.005	2005	
Acatana	0.16	0.0		0.014	0.01	0.038	0.036	0.01
Carbon Disulfide	0.005	0.005		0.005	0.01	0.005	2003	
Xytenes, total	0.005	0.005		0.005	0.005	0.005	0.005	
1,4-Dichlorobenzene				0.01			0.01	
1,2,4-Trichlorobenzene				0.01			0.01	
Bls(2-Ethylhexyl)Phthalate				0.002			0.008	
Methoxychior				0.0005			0.0031	

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	BOTAM	,			BLMM	
	5 [NAS	EPA		EPA	EPA
COMPONICS	Act of	1701/00	201100	ORMOZIL	DAMEN	DAMORA
					,	
2	· ·			-	0.46	
₹		0.01		0.01	0.01	
To the state of th	D. 088	0.177		0.054	0.25	
80					0.005	•
8					0.005	
Ca	43	93,8		52.1	8	
Cr		0.028			0.01	
3 9						
8	ξ.	,			0.014	
2	_				0.025	
Fe		8.37		0.892	0.57	
B	0.0047	0.016			0.003	
160	**************************************	4,16		3.75	Ű٦	
Ma .	0.31	256		0.815	2.2	
2			•		0.04	•
**	₹				Ć1	
Se		0.005		0.005	0.005	
À 0					0.01	
- C		19.8		17.2	÷	-
1					0.01	•
*					0.05	
Zn	0.094				0.065	
Methylene Chloride		0.005	0.013	0.006		0.006
Genzene		0.005	0,013	0.005		0.005
Trichloroethene		0.005	0.013	0.005		2006
Voluena		0.005	0.013	0.005		0.005
Chlerobenzene		0.005	0.013	0.005		6.005
Ethyl Benzene		0.005	0.013	0.005		2000
Acatona		0.148	0.5	0.01		9.
Carbon Disulide		0.005	0.013	0.005		0.005
Xylenes, total		0.005	0.013	0.005		0.005
1,4-Dichlorobenzene			0.01			0.01
1,2,4-Trichloroberzene			0.01			0.01
Bis(2-Ethythexyl)Phthalate			0.008			2002
Methoxychlor			0.0013			0.0005

⊃ 0 N

	MW14					MW14A		
	EPA	SYN	EPA	SYN	EPA .	SYN	SYN	EPA
COMPOUND	1/31/90	1/31/90	2/01/90	4/27/90	7/26/90	1/31/90_	4/27/90	_7/26/90
Ai .	.40				20			4.3
As	0.069	0.078	1	0.029	0.038	0.01	0.01	1
Ba	0.39	0.186	ļ	0.173	0.2	0.393	0.139	0.2
Ве 🐇	0.0024				0.005			0.005
Cd	0.0097				0.005			0.008
Ca	41	64.6		18.9	15	57.9	48.3	49
Cr	0.045	0.017	ĺ	0.013	0.025	0.033	0.01	0.011
Co .	0.036				0.05	į		0.05
Cu	0.067				0.025		` .	0.025
Fe	36	3.72		9.59	11	17	4.03	3.9
Pb	0.09	0.01		0.024	0.025	0.033	0.01	0.0057
Mg	14	2.72		7.25	7.2	4.94	3.12	! 5
Mn /	1.2	0.175		0.376	0.33	3.56	0.695	0.51
Ni	0.13	·	<u> </u>		0.04			0.04
K	76				23	· .		5
Se	0.005	0.007		0.005		0.005	0.005	0.008
Ag	0.01	ł	} '	ł	0.01	ł		0.01
Na	230	146	•	87.9	93	5.66	4.48	5.5
π .	0.01				0.01	ļ	}	0.01
V	0.06		<u> </u>		0.05		1	0.03
2n	0.25	İ			0.074			0.029
Methylene Chloride	Ī	0.05	0.04	0.903	0.005	0.006	0.003	0.005
Benzene		0.05	0.02	0.034	0.005	0.006	0.005	0.005
Trichtoroethène	j	0.05	0.02	0.034	0.005	0.006	0.005	0.005
Toluana	1	0.05	0.02	0.034	0.005	0.006	0.005	0.005
Chlorobanzene	1	0.05	0.02	0.034	0.005	0.006	0.005	0.008
Ehyl Benzene	1	0.05	0.02	0.034	0.005	0.006	0.005	0.005
Acetone		1.34	0.77	0.067	0.5	0.233	0.01	2.1
Carbon Disulikie		9.05	0.2	0.034	0.005	0.008	0.005	0.005
Xylenes, total		0.05	0.02	0.034	0.005	0.006	0.005	0.005
1,4-Dichlorobenzene	1]	0.01]	0.01			0.0
1,2,4-Trichtorobenzene	1	1	0.01		0.01			0.01
Bis(2-Ethylhexyl)Phthalale			0.004	ł	0.01		•	0.0
Methoxychior			0.0005	1				1

TABLE A-3 GROUND WATER DATA ELIMINATED DURING DATA VALIDATION

<u>PARAMETER</u>

Volatiles

Ch loromethane Bromomethane Vinyl Chloride Chloroethane 1,1-Dichloroethene 1,1-Dichloroethane 1.2-Dichloroethens Chloroform 1,2-Dichloroethane 1,1,1-Trichloroethane Carbon Tetrachloride Bromodich loremethane 1,2-Dichloropropane Cis-1.3-Dichloropropene Dibromoch loromethane 1.1,2-Trichloroethane Bronoform Tetrach lorosthene 1,1,2,2-Tetrachloroethane 2-Butone Vinyl Acetate 2-Hexanone 4-Methy 1-2-Pentanone Styrene Trans-1.3-Dichloropropene

Semivolatiles

Trans-1 2-Dichloroethene

2-Chloroethylvinylether

Phonel Bis (2-Chloroethyl) Ether 2-Chlorophenol 1,3-Diahlorobenzene Benzyl Alcohol 1,2-Dichlorobenzen 2-Methy | pheno | Bis (2-Chiloro isopropy 1) Ether 4-Hethy lpheno) N-Nitroso-Dipropylamine Mexach lorgethane Nitrobenzéne 1 sophorene 2-Nitrophenol 2.4-Dimethy iphenol Naphtha lene 4/Chloroaniline Hexach lorobutadiene 4-Chloro-3-Methylphenol 2-Methy inaphtha lens

JUSTIFICATION

Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or apy other sample Not detected in background samples or any other sample Not detected in background samples of any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in Dackground samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not depected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample for detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample

Detected in background samples but not any other sample Not detected in background samples or any other sample Not detected in background samples or any object sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sa Not detected in background samples or any other samp Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Detected in background samples but not any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Detected in background samples but not any other sample



TABLE A-3 GROUND WATER DATA ELIMINATED DURING DATA VALIDATION (CONTINUED)

PARAMETER

JUSTIFICATION

Hexach lorocyc lopentagiene Z. 4.6-Trich larophena? 2,4,5-Trichlorophenol 2-Chloronaphthalene 2-Nitroaniline Dimethylphthalate Acenaphthylene 3-Mitroanilene Acenaphthene 2,4-Dinitrotoluene 2.6-Dinitrotoluene Diethylyghthalate 4-Chlorophenyl Phenyl Ether Fluorene 4-Nitroaniline 4.6-Dinitro-2-Methylphenol N-Nitrosodiphenylamine 4-Bramophenyl Phenyl Ether Hexach larobenzene Pentach loropheno l Phenanthrene Anthracese Di-N-Buty lPhtha late Fluoranthene Pyrene Butyl Benzyl Phthalate 3.3-Dichlorobenzidine Benzo (A) Anthracene Chrysene Di-N-Octyl Phthalate Benzo(B)Fluoranthene Benzo (K) Fluoranthene Indeno(1.2.3-CD)Pyrene Dibenzo (A.H) Anthracene

Pesticides
Alpha-BHC
Beta-BHC

Benzo (G. H<1) Pery lene

Delta-BHC

Gamma-BHC

Aldrin Diejdrin

A Endozulfan (1)

B/Endosulfan (II) Ændosulfan Sulfate

Endrin

Endrin Aldehyde

4,4-00E 4,4-00D

Not detected in background samples or app other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in packground samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in Background samples or any other sample Not detegred in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or and other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other\sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample

Not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample not detected in background samples or any other sample

TABLE A-3 GROUND WATER DATA ELIMINATED DURING DATA VALIDATION (CONTINUED)

JUSTIFICATION

' \	
4.4°DDT	Not detected in background samples or any other sample
Toxaphene	Not detected in background samples or shy other sample
PCB-1016	Not detected in background samples or any other sample
PCB-1221	Not detected in background samples or any other sample
PCB-1232	Not detected in background samples or any other sample
PC8-1242	Not detected in background samples or any other sample
PCB-1248	Not detected in background samples or any other sample
PCB-1254	Not detected in background samples or any other sample
PCB-1280	Not detected in background samples or any other sample
Chilorodane Technical	Not detected in background samples or any other sample
Heptachlor	Not detected in background samples or any other sample
Heptach for Epoxida	Not detected in background samples or any other sample
Endrin Ketone	Not detected in background samples or any other sample
Alpha-Chlurdase	Not detected in background samples or any other sample
Gamia-Chirdane	Not detected in background samples or any other sample

Total Metals

Antimony Moreury Cyanide

ted in background samples or any other sample Detected in background samples but not any other sample nalyzed in any sample

Dissolved Metals

Aluminum Ant imony Beryllium Chromium Copper Nickel Selentum Silver Tha 11 tun Vanad ium Cyanide

Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not detected in background samples or any other sample Not analyzed in any sample

Individual Data Points

MW14 EPA Split Sample dated 7/26/90

Detection limit is higher than the Contract Required Detection Limit Se len ium

HW 1 EPA Split/ Sample dated 7/26/90

Detection limit is higher than the Contract Required Detection Limit

MW5 EPA Split Sample dated 7/26/90

Detection limit is higher than the Contract Required Detection Limit Thallium.

MV6 EPA Split Sample dated 7/26/90

Se len tuni Detection limit is higher than the Contract Required Detection Limit Tha llium Detection limit is higher than the Contract Required Detection Limit

HW158/EPA Split Sample dated 7/26/90

The 1/1 um Detection limit is higher than the Contract Required Detection Limit

Mw16B EPA Split Sample dated 7/26/90

Se/lenium Detection limit is higher than the Contract Required Detection Limit

TABLE A-3 GROUND WATER DATA ELIMINATED DURING DATA VALIDATION (CONTINUED)

NOTES:

MW5 EPA Split Sample dated 7/26/90 had an unusally high value for total lead. This value falls out of the 95% confidence interval by a significant margin. This data is qualified by the laboratory, however excluding the data from this set is not justified. A resampling event may be necessary to verify this data.

	SYNTEX					BACKGRO	UND		
				STANDARD	95% C.I.				REASON FOR ELIMINATION FROM
COMPOUND	MINIMUM	MAXIMUMI	MEAN	DEVIATION		MINIMUM	MAXIMUM	MEAN	CONSIDERATION AS CHEMICAL OF CONCER
Al	9.2	EUV .	75.084	154,50198	384.088405	4.3	40	21.43333	TOXICITY DATA INADEQUATE
As	0.01		0.0164		0.04430854	0.01		0.039	BACKGROUND HIGHER
Ba ·	0.047		1.6821	2.2068380	6.09583739	0.139		0.240142	BASKSHOOND HEILEN
Be	0.0028		6:Q114		0.04126760	0.0024		0.004133	
Cd ·	0.0024			0.0195844	0.05242606	0.005		0.806566	
Ca	0.0024		98.117	\	187.660034	15	64.6		ESSENTIAL NUTRIENT
Or Or	0.005		1.4681	5.7818626	13.0356259	0.01	0.045	0.022	COSCA FIAC NOT HILLY
Co .	0.0081		0.0722	0.0865940	0.24542137	0.036	0.045		ENZYME COFACTORINUTRIENT
Cu	0.025	* - *	0.0847	0.1138054	0.31236096	0.035	0.067	0.039	ESSENTIAL NUTRIENT
Fa	0.023		84.453	152.49499	389.443720	3.72	0.057 36/	12.17714	ESSENTIAL NUTRIENT
ru Po	0.001		7.9942	41.955379	91.9050241	8-Q057	_0.09	0.028242	LEGICITIES INC. INC. INC. INC. INC. INC. INC. INC.
ro Vig	2,41		11.269	7.6052481	26,4798511	2.78	14		LEVELS SIMILAR TO BACKGROUND
vin .	0.204		12.344	19.309955	50.9643946	9:175	3.58		ESSENTIAL NUTRIENT
Al.	0.204		0.1053		0.83017173	0.04	0.30	0.570071	ESSERTIAL ING TRICAT
K	5	53	1255	14.209028	40.9680576	5	76		BACKGROUND HIGHER
N Se	0.005		0.0084	0.0119911	0.03244378	0.005	0.007		ESSENTIAL NUTRIENT
	0.003		0.0063		0.03858116	0.005	0.007	•	I.
Ag Na	0.01	••••			127.985662	4.46	230		BACKGROUND HIGHER
1981. Ti		0.1		0.0186125	0.04222518	0.01	0.01		LEVELS SIMILAR TO BACKGROUND
V	0.01			0.1407451	0.32471617	0.05			LEVELS SIMILAR TO BACKGROUND
v Zn	0.028		/	7.1198822	16,2182160	0.029	0.05	0.053333	LEVELS STRILAN TO BACKGROUND
	0.055		1.9784		0.30962603	0.029	0.903		BACKGROUND HIGHER
Asthylene Chloride	0.001	<i>_</i> ·	0.0256	0.0617270	0.30952603	0.005	0.903	0.017857	BUNGHOOM HOUCK
Jenzene	0.001					1		0.017857	
(richioroethene	0.005	0.25 0.25	0.026	0.081 0 127 0.0662629	0.14922540 0.16355717	0.005 0.005	0.05 0.05	0.017857	
Toluene	0.005					1		0.017857	
Chlorobenzene	0.002		0.0416	0.0670292	0.17568352	0.005	0.05 0.05		
Ethyl Benzene	0.005	-	0.0271		0.14995904	0.005		0.017857	COMMON I ADODATODY CONTANTANT
Cetone	6.01		1.1843	2.6783542	8.54105222	0.01	21		COMMON LABORATORY CONTAMINANT
Carbon Disulfide	0.002		0.0258	0.0610686	0.14914988	0.005	0.2	0.043571	LEVELS SIMILAR TO BACKGROUND
(yienes, total	0.001		0.0360	0.0657677	0.16762925	0.005		0.017857	
A-Dichlorobenzene	0.01	-	0.0228	0.0293158	0.08144996	0.01	0.01	0.01	LEGIC CONTRACTOR OF CASCASTA
2,4-Trichlorobenzens	0.01		0.0100	0.0002874	0.01066586	0.01	0.01		LEVELS SIMILAR TO BACKGROUND
Bis(2-Ethylhexyl)phthlate	0.002	0.015	0.008	0.0038847	0.01576940	0.004	0.01	0.008	LEVELS SIMILAR TO BACKGROUND
Methoxychlor	0.0005	0.0031	0.0012	0.0006892	0.00307352	0.0005	0.0005	0.0005	

& XIQN399A

Toxicity Profiles

(to be provided)

TABLE TOXICITY FACTORS FOR QUANTIFICATION OF CHRONIC AND LIFETIME HAZARDS FOR CONSTITUENTS OF CONCERN (1)

	CHRONI		CARCINOGEN SLO	PE FACTORS	
	NONCARCINOG	EN RIDS	CANCER	CANCER	
CHEMICAL.	ORAL RID (MG/KG/DAY)	INHALATION RID (MG/KG/DAY)	SLOPE FACTOR (ORAL) (MG/KG/DAY)-1	SLOPE FACTOR (INHALATION) (MG/KG/DAY)-1	CANCER WEIGHT OF EVIDENCE
AL UMINUM	NA	NA	NA	NA	Α .
BARIUM	7.00E-02	NA	. NA	NA	NA:
BERYLLIUM	5.00E-03	NA .	4.30E+00	8.40E+00	B2
<u>CADMUN</u>	5.00E-04	NA	NA	6.10E+09	B1
CHROMIUM (2)	5.00E-03	NA	NA	4.10E+01	· A
LEAD	NA	NA	NA	NA	B2
NICKE	2.00E-02	NA	. NA	NA -	NA
ZINC (3)	2.00E-01	NA	NA	NA	NA
BENZENE	MA	NA	2.90E-02	2,90E-02	A
TRICHLOROETHENE (4)	MA	NA .	1.10E-02	1.70E-02	B2 ·
TOLUENE	2.00E-01	5.70E-01 (5)	NA ·	NA	D
CHLOROBENZENE	2.00E-02	5.00E-03 (6)	NA	NA	D
ETHYL BENZENE	1.00E-01	NA"	NA	NA	D
XYLENES, total	2.00E+00	3.00E-01 (6)	NA	NA	D
1,4-DICHLOROBENZENE	NA	7.00E-01 (6)	2.40E-02 (4)	NA	92
METHOXYCHLOR	5.00E-03	NA	NA	NA	D

- (1) Source: IPIS, November 27, 1990 (unless otherwise indicated).
- (2) The RID and Cancer Slope Factor for chromium reflects the most conservative value provided for either chromium (II) chromium (VI): Noncarcinogen Oral Rid chromium (VI)

 Inhalation Cancer Slope Factor chromium (VI)
- (3) Source for zinc oral RfD values: U.S. EPA, 1989b.
- (4) Source for trichloroethene/1,4-dichlorobenzene CSF values: U.S. EPA, 1989b.
- (5) Source for toluene RID values: U.S. EPA, 1989b. Value derived assuming 70 kg body weight and 20 m3/day inhalation rate.
- (6) Source for chlorobenzene/xylenes/1-4 dichlorobenzene inhalation RID values: U.S. EPA, 1989b.

TABLE TOXICITY FACTORS FOR QUANTIFICATION OF CHRONIC AND LIFETIME HAZARDS FOR CONSTITUENTS OF CONCERN (1)

CHRONIC

CARCINOGEN SLOPE FACTORS

	NONCARCINOGEN RIDE	EN PIDS	OANCED	CANCED
		INHALATION	SLOPE FACTOR	SLOPE FACTOR
CHENICAL	(MGIKG/DAY)	RID (MG <i>INGJID</i> AY)	(ORAL) MG/NG/DAY)-1	(MICKG/DAY)-1
BARIUM	7.00E-02	AA	NA	AN
BERYLLIUM	5.00E-03	*	4.30E+00	8.40m-00
CADMIUM	5.00E-04	¥	×	8.10E - 00
CHROMIUM (2)	5,00E-03	*	NA.	4,10E+01
LEAD	NA A	×	NA	*
NICHEL	2.00E-02	*	NA.	NA.
ZINC(3)	2.00€-01	NA	NA	N _P
BENZENE	· *	NA.	2,90E-02	2.90E-02
TRICHLOROETHENE (4)	NA	K)	1, 10E-02	1.70E-02
TOLUENE	2.00E-01	5.70E-01 (5)	NA.	3
CHLOROBENZENE	2.00E-02	•	N.	¥ A
ETHYL BENZENE	1.00E-01		N,	NA.
XYLENES, total	2.00E+00	3.00E-01 (6)	N.	NA NA
1,4-DICHLOHOBENZENE	₹	_	2.405-02 (4)	NA
METHOXYCHLOR	5.00E-03	Z>	2×	RA

chromium (VI): Norcascinogen Oral Rtd - chromium (VI)

多数多克 Source for zinc oral RID values: U.S. EPA, 1989b.

Inhalation Cancer Slope Factor - chromium (VI)

(1) Source: IRIS, November 27, 1990 (unless otherwise indicated).
(2) The RfD and Cancer Slope Factor for chromium reflects the most conservative value provided for either chromium (III)

Source for trichloroethens/1,4-dichlorobenzene CSF values: U.S. EPA, 1989th.

Source for toluene RID values: U.S. EPA, 1989b. Value derived assuming 70 kg body weight and 20 m3/day inhalation rate.

Source for chlorobenzene/xyfenes/1-4 dichlorobenzene Inhalation RID values: U.S. EPA, 1989b.

TABLE TOXICITY FACTORS FOR QUANTIFICATION OF SUBCHRONIC HAZARDS FOR CONSTITUENTS OF CONCERN (1)

SUBCHRONIC RIDS FOR MONCARCINOGENS.

Y Baic All	(MOJKG/DAY)	(MORG/DAY)
8A.P.L.M.	5.00E-02	1.00E-03
	5.00E-00	ND
Xoxius	CS	2
HOMUM (2)	2.00E-02	S
503	S	ND
SOXEL	2,00€-02	S
	2,00€-01	8
	NO	20
RICH LOROETHENE	₹B	NO O
	4.00E-01	5.70E-01
HOROBENZENE	2.00E-01	5.00E-03
ETYLBENZENE	1.00€+00	2
O'LENES local	4.00€+00	8.57E-02
,4-DICHLOROBENZENE	9.00E-01	4.00E-01
METHOXYCHLOR	1.00E-01	2

⁽³⁾ Final draft all quality criteria document (600\0-63-028F) declines to derive an air quality criterion for isad.
(4) Value dived assuming 70 kg body weight and 20 m3/day inhalation rate.

DERMAL HAZARDS AND RISKS ASSOCIATED WITH INDICATOR CHEMICALS

	ABCOSOTION	VALUES VALUES	JES JES	VAI	VALUES	CARCIN CODA ANTECES	DA AVERSES
CHEMICAL	FACTORS USED FOR ORAL TOXICITY VALUES (2)	ADJUSTED(2 SUBCHRONIC SUBCHRONI ORAL HID RID (MG/KG/DAY) (MG/KG/DAY)	ADJUSTEX(2) SUBCHRONIC SUBCHRONIC ORAL RID RID (MG/KG/DAY) (MG/KG/DAY)	CHRONIC ORAL RID (MG/KG/DAY)	ADJUSTED(3) CHRONIC RIC (MG/KG/DAY) (CANCER SLOPE FACTOR (ORAL) MGMG/DAY)-1	ADJUSTED(4) SLOPE (MG/KG/DAY)-1
BARIUM	10.00	5.00E-02	5.00E-03	7.00E-02	7.00E-03	NA	NA.
BERYLLIUM	10.00		5.00E-04	5.00€-03	5.00E-04	4.30E+00	4.30E+01
CADMIUM	7.00		N.A.	5.00E-04		Z	Z P
CHROMIUM (2)	50,00	200E-02	1,00E-02	5,00E-03		NA A	NA
.EAD (3)	50,00	Z.	N A	Z	N,	Z A	ş
NICKEL	10.00	2.00E-02	2.00E-03	200E-02		N A	≩
ZINC	30.00		6.00€-02	200E-01	6.00E-02	Ş	¥
HENZENE	90.00		¥	ž	Z	2.90E-02	3.22E-02
FRICHLORIDETHENE	100.00	Z	NA A	×	Z	1.10E-02	1.10E-02
COLUENE	100.00	4.00E-01	4.00E-01	200E-01	2.00E-01	N _A	Z ≯
CHLOROBENZENE	30.00		6,00E-02	200E-02	6.00E-	NA A	₹
ETHYLBENZENE	90.00		8.00E-01	1.00E-01	8.00E-02	NA	₹
XYLENES, TOTAL	92.00	4.00E+00	3.68€+00	200E+00	1.84E+00	NA A	₹
,4-DICHLOFIOBENZENE	100.00	9.00E-01	9.00E-D1	Z *	0.00E+00	2,40E-02	2.40E-02
METHOXYCHLOR	100.00	1.00E-01	1.00E-D1	5.000-03	5.00E-03	\$	\$

Sources for subchronic and chronic RID/CSF values = U.S. EPA, 1990; U.S. EPA, 1988; ECAO, 1994; U.S. EPA, 1988b.
 Oral absorbtion factors are discussed in Appendix D (Source U.S.EPA, 1990).
 Adjustment of an administered to an absorbed dose RfD:

(7) Because PAHs such as benzo(a)pyrene are skin carcinogens which act directly at the point of contact, it is inappropriate to use the craft slope factor to exalt ate dermal cancer risks for these compounds.

TIVE

⁽Administered RID) x (Oral Absorption Factor) = Absorbed Dose RID

Ī Adjustment of an administered to an absorbed dose stope factor.

⁽Administered Slope Factor)-1 / (Oral Absorption Factor)

⁽⁶⁾ Toxicity factors for PAHs are based on values for benzo(a) pyrene. (5) The RID and Cancer Slope Factor for chromium reflects the most conservative value provided for either chromium (III) 8 chromium (VI): Noncarcinogen Oral Rfd - chromium (VI)